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UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION 10 LABORATORY  
7411 BEACH DR. EAST  
PORT ORCHARD, WASHINGTON 98366

October 13, 1992

MEMORANDUM

SUBJECT: Report of Data Validation of BNA's for the Avery Railroad Project, Water Samples 92352351, 92352352, 92352353 and 92352354.

FROM: J. Blazeovich, Program Manager  
Analytical Chemistry Program *JNB*

TO: Monica Rolluda, Project Officer  
Avery Railroad Project

CC: *BRUCE WOODS*  
*KARA STEWARD*

The following is a QA data review of the BNA analysis of water samples collected for the Avery Railroad project and performed at the Manchester Laboratory. This review covers the following samples:

92352351 92352352 92352353 92352354

The project code for these samples is TEC-575A and the account number is 2TFA10PUZZ.

Data qualifications

The following comments refer to laboratory performance meeting the Quality Control specifications outlined in the "CLP Statement of Work, Organic Analysis, revision 1/91."

I. Holding Times: Acceptable

All samples were held seven days or less between collection and extraction. All extracts were held less than forty days between extraction and analysis. No data qualifiers were required due to exceeding holding times.

II. GC/MS Tuning and Performance: Acceptable

The tuning summary agreed with the raw data. All deca-fluorotriphenylphosphine ion abundance met criteria. All sample analyses were preceded by a tune less than 12 hours prior to analysis. No data qualifiers were required on the basis of tuning data.

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### III. Initial Calibration: Acceptable

An extended seven point calibration curve was constructed for most target analytes on 1/17/92. The calculation of the relative response factors was checked and the calculation method was correct. All analytes met the SPCC criterion (0.05). The %RSD exceeded 30% for six targets, N-nitrosodiphenylamine, benzoic acid, 4,6-dinitro-2-methylphenol, 2,4-dinitrophenol, carbazole and di-n-octyl phthalate. Each were given the data qualifier J or UJ. All other response factors and %RSDs were acceptable.

### IV. Continuing Calibration: Acceptable

9/28/92 Sample 92352351, 92352352, 92352353, 92352354, Matrix Spikes 92352353Y, 92352353Z and Blanks EBW2245D1, EBW2245D2.

The response factors for all compounds were above 0.05 except 3-nitroaniline a had response factor of 0.032. All negative results for this analyte were qualified as unusable, R. Eight compounds, benzyl alcohol, 2,6-dinitrotoluene, 4-chloro-3-methylphenol, diethyl phthalate, di-n-butyl phthalate, carbazole butylbenzyl phthalate, and bis(2-ethylhexyl) phthalate, had %diff. values greater than 30%. All values for these analytes were given the qualifier J or UJ. No additional qualifiers were required on the basis of response factor or %diff.

### V. Blanks: Acceptable

Two blanks were analyzed with the samples. Common laboratory contaminants, N-nitrosodiphenylamine and bis(2-ethylhexyl) phthalate along with low concentrations of naphthalene, 2-methylnaphthalene, 1-methylnaphthalene, acenaphthalene and phenanthrene, were detected. These analytes were reported in the samples only if the associated integrated areas exceeded those found in the corresponding blank by ten times. No data qualifiers were required based on blank results.

### VI. Surrogates: Acceptable

All surrogate recoveries were within specification for all blanks, spikes and samples except for the diluted sample extract of 92352354. Since the extract was diluted, some of the surrogates were not detected. No data qualifiers were required.

### VII. Matrix Spike/Matrix Spike Duplicate: Acceptable

Recoveries of most MS/MSD analytes were within CLP and/or Region Ten guidelines. One or both spike recoveries for four analytes, hexachloroethane, hexachlorocyclopentadiene, benzoic acid and 3-nitroaniline, were outside the 50-150% Region Ten acceptance window. All values for these analytes were given the qualifier J or UJ for the corresponding sample, 92352353. No

additional data qualifiers were required on the basis of MS/MSD results.

VIII. Internal Standard Performance: Acceptable

The retention time variations of all internal standards were within 30 seconds of the daily standard which is acceptable. The %area of all the internal standards fell within the 50-200% acceptance window. No data qualifiers were required on the basis of internal standards data.

IX. TCL Compound Identification: Acceptable

All detected TCL compounds' relative retention times were within acceptable limits of the related standards in the continuing calibration standard. Criteria were met for mass spectral ion matching and ion abundance matching or the mass spectra were judged acceptable.

X. Compound Quantitation: Acceptable

Compound quantitation was evaluated correctly. The appropriate internal standards were used. The correct quantitation ions and relative response factors were used. Some analytes were detected in this sample set at levels below the lowest calibration concentration of the initial calibration curve. Since the Practical Quantitation Limit is based on this lowest initial calibration standard, these values were assigned the J qualifier. No additional qualifiers were required on the basis of compound quantitation.

XI. Tentatively Identified Compounds: Acceptable

Spectra for all tentatively identified compounds met criteria for mass spectral ion matching and ion abundance matching or the mass spectra were judged acceptable. Some unknown compounds were detected.

Overall Assessment for the Case

The usefulness of the data is based on the criteria outlined in the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses (6/91)."

All requirements for data qualifiers from the preceding sections were accumulated. Each sample data summary sheet and each compound were checked for positive or negative results. From this the overall need for data qualifiers for each analysis was determined. In cases where more than one of the preceding sections required data qualifiers, the most restrictive qualifier has been added to the data.

In general, all unqualified data can be used without restriction. The usefulness of qualified data should be treated according to the severity of the qualifier. Should questions arise regarding the qualification of data and its relation to the usefulness, the reader is encouraged to contact the Region 10 laboratory.

#### DATA QUALIFIERS

- U - The analyte was not detected at or above the reported result.
- J - The analyte was positively identified. The associated numerical result is an estimate.
- EXP - The result is equal to the number before EXP times 10 to the power of the number after EXP. As an example 3EXP6 equals  $3 \times 10^6$ .
- REJ - The data are unusable for all purposes.
- N - There is evidence the analyte is present in this sample.
- NJ - There is evidence that the analyte is present. The associated numerical result is an estimate.
- UJ - The analyte was not detected at or above the reported estimated result. The associated numerical value is an estimate of the quantitation limit of the analyte in this sample.
- NAF - Not analyzed for.
- \* - The analyte was present in the sample.  
(Visual aid to locate detected compounds on the report sheet.)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Laboratory: EPA, Manchester

Sample No: 92 352351

Description: WAW01

Source: Well (Drinking Water Supply)

Begin Date: 92/08/26 08:37

B/N/Acid Scan			B/N/Acid Scan		
Water-Total		Units	Water-Total		Units
Result			Result		
Benzo(a)pyrene	2U	ug/l	Di-n-Octyl Phthalate	2UJ	ug/l
2,4-Dinitrophenol	26UJ	ug/l	Hexachlorobenzene	2U	ug/l
Dibenzo(a,h)anthracene	5U	ug/l	Anthracene	2U	ug/l
Benzo(a)anthracene	2U	ug/l	1,2,4-Trichlorobenzene	2U	ug/l
4-Chloro-3-Methylphenol	10UJ	ug/l	2,4-Dichlorophenol	2U	ug/l
Benzoic acid	26UJ	ug/l	2,4-Dinitrotoluene	5UJ	ug/l
Hexachloroethane	2U	ug/l	Pyrene	2U	ug/l
Hexachlorocyclopentadi+	10U	ug/l	Dimethylphthalate	2U	ug/l
Isophorone	2U	ug/l	Dibenzofuran	2U	ug/l
Acenaphthene	2U	ug/l	Benzo(ghi)perylene	2U	ug/l
Diethylphthalate	2UJ	ug/l	Indeno(1,2,3-cd)pyrene	2U	ug/l
Di-n-Butylphthalate	2UJ	ug/l	Benzo(b)fluoranthene	2U	ug/l
Phenanthrene	2U	ug/l	Fluoranthene	2U	ug/l
Butylbenzylphthalate	5UJ	ug/l	Benzo(k)fluoranthene	2U	ug/l
N-Nitrosodiphenylamine	26UJ	ug/l	Acenaphthylene	2U	ug/l
Fluorene	2U	ug/l	Chrysene	2U	ug/l
Carbazole	10UJ	ug/l	Retene	2U	ug/l
Hexachlorobutadiene	5U	ug/l	4,6-Dinitro-2-methylph+	26UJ	ug/l
Pentachlorophenol	10U	ug/l	1,3-Dichlorobenzene	2U	ug/l
2,4,6-Trichlorophenol	5U	ug/l	2,6-Dinitrotoluene	5U	ug/l
2-Nitroaniline	5U	ug/l	N-Nitroso-di-n-Propyla+	2U	ug/l
2-Nitrophenol	5U	ug/l	4-Chlorophenyl-phenyle+	2U	ug/l
Naphthalene, 1-Methyl-	2U	ug/l	bis(2-Chloroisopropyl)+	2U	ug/l
Naphthalene	2U	ug/l	Surrog: 2-Fluorobiphen+	80	% Recov
2-Methylnaphthalene	2U	ug/l	Surrog: 2-Fluorophenol	68	% Recov
2-Chloronaphthalene	2U	ug/l	D4-1,2-Dichlorobenzene	50	% Recov
3,3'-Dichlorobenzidine	52U	ug/l	Surrog: D14-Terphenyl	68	% Recov
2-Methylphenol	2U	ug/l	PYRENE-D10 (SS)	71	% Recov
1,2-Dichlorobenzene	2U	ug/l	Surrog: D5-Nitrobenzene	74	% Recov
o-Chlorophenol (2-Chlo+	2U	ug/l	Surrog: D5-Phenol	56	% Recov
2,4,5-Trichlorophenol	10U	ug/l			
Nitrobenzene	2U	ug/l			
3-Nitroaniline	REJ	ug/l			
4-Nitroaniline	26U	ug/l			
4-Nitrophenol	13U	ug/l			
Benzyl Alcohol	41UJ	ug/l			
4-Bromophenyl-phenylet+	2U	ug/l			
2,4-Dimethylphenol	2U	ug/l			
4-Methylphenol	2U	ug/l			
1,4-Dichlorobenzene	2U	ug/l			
4-Chloroaniline	26U	ug/l			
Phenol	2U	ug/l			
bis(2-Chloroethyl)Ether	2U	ug/l			
bis(2-Chloroethoxy)Met+	2U	ug/l			
BIS(2-ETHYLHEXYL) PHTH+	2UJ	ug/l			

(Sample Complete)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Laboratory: EPA, Manchester

Sample No: 92 352352

Description: WSW01

Source: Well (Drinking Water Supply)

Begin Date: 92/08/26 09:15

B/N/Acid Scan			B/N/Acid Scan		
Water-Total		Units	Water-Total		Units
Result			Result		
Benzo(a)pyrene	2U	ug/l	Di-n-Octyl Phthalate	2UJ	ug/l
2,4-Dinitrophenol	27UJ	ug/l	Hexachlorobenzene	2U	ug/l
Dibenzo(a,h)anthracene	5U	ug/l	Anthracene	2U	ug/l
Benzo(a)anthracene	2U	ug/l	1,2,4-Trichlorobenzene	2U	ug/l
4-Chloro-3-Methylphenol	11UJ	ug/l	2,4-Dichlorophenol	2U	ug/l
Benzoic acid	27UJ	ug/l	2,4-Dinitrotoluene	5UJ	ug/l
Hexachloroethane	2U	ug/l	Pyrene	2U	ug/l
Hexachlorocyclopentadi+	11U	ug/l	Dimethylphthalate	2U	ug/l
Isophorone	2U	ug/l	Dibenzofuran	2U	ug/l
Acenaphthene	2U	ug/l	Benzo(ghi)perylene	2U	ug/l
Diethylphthalate	2UJ	ug/l	Indeno(1,2,3-cd)pyrene	2U	ug/l
Di-n-Butylphthalate	2UJ	ug/l	Benzo(b)fluoranthene	2U	ug/l
Phenanthrene	2U	ug/l	Fluoranthene	2U	ug/l
Butylbenzylphthalate	5UJ	ug/l	Benzo(k)fluoranthene	2U	ug/l
N-Nitrosodiphenylamine	27UJ	ug/l	Acenaphthylene	2U	ug/l
Fluorene	2U	ug/l	Chrysene	2U	ug/l
Carbazole	11UJ	ug/l	Retene	2U	ug/l
Hexachlorobutadiene	5U	ug/l	4,6-Dinitro-2-methylph+	27UJ	ug/l
Pentachlorophenol	11U	ug/l	1,3-Dichlorobenzene	2U	ug/l
2,4,6-Trichlorophenol	5U	ug/l	2,6-Dinitrotoluene	5U	ug/l
2-Nitroaniline	5U	ug/l	N-Nitroso-di-n-Propyla+	2U	ug/l
2-Nitrophenol	5U	ug/l	4-Chlorophenyl-phenyle+	2U	ug/l
Naphthalene, 1-Methyl-	2U	ug/l	bis(2-Chloroisopropyl)+	2U	ug/l
Naphthalene	2U	ug/l	Surrog: 2-Fluorobiphen+	85	% Recov
2-Methylnaphthalene	2U	ug/l	Surrog: 2-Fluorophenol	74	% Recov
2-Chloronaphthalene	2U	ug/l	D4-1,2-Dichlorobenzene	53	% Recov
3,3'-Dichlorobenzidine	54U	ug/l	Surrog: D14-Terphenyl	76	% Recov
2-Methylphenol	2U	ug/l	PYRENE-D10 (SS)	80	% Recov
1,2-Dichlorobenzene	2U	ug/l	Surrog: D5-Nitrobenzene	81	% Recov
o-Chlorophenol (2-Chlo+	2U	ug/l	Surrog: D5-Phenol	62	% Recov
2,4,5-Trichlorophenol	11U	ug/l			
Nitrobenzene	2U	ug/l			
3-Nitroaniline	REJ	ug/l			
4-Nitroaniline	27U	ug/l			
4-Nitrophenol	13U	ug/l			
Benzyl Alcohol	43UJ	ug/l			
4-Bromophenyl-phenylet+	2U	ug/l			
2,4-Dimethylphenol	2U	ug/l			
4-Methylphenol	2U	ug/l			
1,4-Dichlorobenzene	2U	ug/l			
4-Chloroaniline	27U	ug/l			
Phenol	2U	ug/l			
bis(2-Chloroethyl)Ether	2U	ug/l			
bis(2-Chloroethoxy)Met+	2U	ug/l			
BIS(2-ETHYLHEXYL) PHTH+	2UJ	ug/l			

(Sample Complete)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Laboratory: EPA, Manchester

Sample No: 92 352353

Description: WSW02

Source: Well (Drinking Water Supply)

Begin Date: 92/08/26 09:15

B/N/Acid Scan			B/N/Acid Scan			B/N/Acid Scan		
Water-Total			Water-Total			Water-Total		
Result Units			Result Units			Result Units		
*** Continued ***			*** Continued ***			*** Continued ***		
Matrix Spike #1			Matrix Spike #1			Matrix Spike #1		
Benzo(a)pyrene	2U	ug/l	Di-n-Octyl Phthalate	2UJ	ug/l	Acenaphthene	74	% Recov
2,4-Dinitrophenol	25UJ	ug/l	Hexachlorobenzene	2U	ug/l	Diethylphthalate	81J	% Recov
Dibenzo(a,h)anthracene	5U	ug/l	Anthracene	2U	ug/l	Di-n-Butylphthalate	75J	% Recov
Benzo(a)anthracene	2U	ug/l	1,2,4-Trichlorobenzene	2U	ug/l	Phenanthrene	78	% Recov
4-Chloro-3-Methylphenol	10UJ	ug/l	2,4-Dichlorophenol	2U	ug/l	Butylbenzylphthalate	67J	% Recov
Benzoic acid	25UJ	ug/l	2,4-Dinitrotoluene	5UJ	ug/l	N-Nitrosodiphenylamine	124J	% Recov
Hexachloroethane	2UJ	ug/l	Pyrene	2U	ug/l	Fluorene	78	% Recov
Hexachlorocyclopentadi+	10UJ	ug/l	Dimethylphthalate	2U	ug/l	Carbazole	NAF	% Recov
Isophorone	2U	ug/l	Dibenzofuran	2U	ug/l	Hexachlorobutadiene	53	% Recov
Acenaphthene	2U	ug/l	Benzo(ghi)perylene	2U	ug/l	Pentachlorophenol	71	% Recov
Diethylphthalate	2UJ	ug/l	Indeno(1,2,3-cd)pyrene	2U	ug/l	2,4,6-Trichlorophenol	66	% Recov
Di-n-Butylphthalate	2UJ	ug/l	Benzo(b)fluoranthene	2U	ug/l	2-Nitroaniline	82	% Recov
Phenanthrene	2U	ug/l	Fluoranthene	2U	ug/l	2-Nitrophenol	90	% Recov
Butylbenzylphthalate	5UJ	ug/l	Benzo(k)fluoranthene	2U	ug/l	Naphthalene, 1-Methyl-	NAF	% Recov
N-Nitrosodiphenylamine	25UJ	ug/l	Acenaphthylene	2U	ug/l	Naphthalene	70	% Recov
Fluorene	2U	ug/l	Chrysene	2U	ug/l	2-Methylnaphthalene	52	% Recov
Carbazole	10UJ	ug/l	Retene	2U	ug/l	2-Chloronaphthalene	70	% Recov
Hexachlorobutadiene	5U	ug/l	4,6-Dinitro-2-methylph+	25UJ	ug/l	3,3'-Dichlorobenzidine	NAF	% Recov
Pentachlorophenol	10U	ug/l	1,3-Dichlorobenzene	2U	ug/l	2-Methylphenol	74	% Recov
2,4,6-Trichlorophenol	5U	ug/l	2,6-Dinitrotoluene	5U	ug/l	1,2-Dichlorobenzene	60	% Recov
2-Nitroaniline	5U	ug/l	N-Nitroso-di-n-Propyla+	2U	ug/l	o-Chlorophenol (2-Chlo+	80	% Recov
2-Nitrophenol	5U	ug/l	4-Chlorophenyl-phenyle+	2U	ug/l	2,4,5-Trichlorophenol	77	% Recov
Naphthalene, 1-Methyl-	2U	ug/l	bis(2-Chloroisopropyl)+	2U	ug/l	Nitrobenzene	76	% Recov
Naphthalene	2U	ug/l	Surrog: 2-Fluorobiphen+	89	% Recov	3-Nitroaniline	REJ	% Recov
2-Methylnaphthalene	2U	ug/l	Surrog: 2-Fluorophenol	74	% Recov	4-Nitroaniline	81	% Recov
2-Chloronaphthalene	2U	ug/l	D4-1,2-Dichlorobenzene	56	% Recov	4-Nitrophenol	58	% Recov
3,3'-Dichlorobenzidine	50U	ug/l	Surrog: D14-Terphenyl	84	% Recov	Benzyl Alcohol	66J	% Recov
2-Methylphenol	2U	ug/l	PYRENE-D10 (SS)	86	% Recov	4-Bromophenyl-phenylet+	72	% Recov
1,2-Dichlorobenzene	2U	ug/l	Surrog: D5-Nitrobenzene	84	% Recov	2,4-Dimethylphenol	68	% Recov
o-Chlorophenol (2-Chlo+	2U	ug/l	Surrog: D5-Phenol	58	% Recov	4-Methylphenol	74	% Recov
2,4,5-Trichlorophenol	10U	ug/l				1,4-Dichlorobenzene	59	% Recov
Nitrobenzene	2U	ug/l				4-Chloroaniline	110	% Recov
3-Nitroaniline	REJ	ug/l				Phenol	61	% Recov
4-Nitroaniline	25U	ug/l				bis(2-Chloroethyl)Ether	80	% Recov
4-Nitrophenol	12U	ug/l				bis(2-Chloroethoxy)Met+	75	% Recov
Benzyl Alcohol	40UJ	ug/l				BIS(2-ETHYLHEXYL) PHTH+	63J	% Recov
4-Bromophenyl-phenylet+	2U	ug/l	Benzo(a)pyrene	73	% Recov	Di-n-Octyl Phthalate	62J	% Recov
2,4-Dimethylphenol	2U	ug/l	2,4-Dinitrophenol	64J	% Recov	Hexachlorobenzene	75	% Recov
4-Methylphenol	2U	ug/l	Dibenzo(a,h)anthracene	71	% Recov	Anthracene	79	% Recov
1,4-Dichlorobenzene	2U	ug/l	Benzo(a)anthracene	74	% Recov	1,2,4-Trichlorobenzene	59	% Recov
4-Chloroaniline	25U	ug/l	4-Chloro-3-Methylphenol	69J	% Recov	2,4-Dichlorophenol	67	% Recov
Phenol	2U	ug/l	Benzoic acid	5J	% Recov	2,4-Dinitrotoluene	78J	% Recov
bis(2-Chloroethyl)Ether	2U	ug/l	Hexachloroethane	48	% Recov	Pyrene	78	% Recov
bis(2-Chloroethoxy)Met+	2U	ug/l	Hexachlorocyclopentadi+	20	% Recov	Dimethylphthalate	80	% Recov
BIS(2-ETHYLHEXYL) PHTH+	2UJ	ug/l	Isophorone	72	% Recov			

(Continued on next page)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Laboratory: EPA, Manchester

Sample No: 92 352353

Description: WSW02

Source: Well (Drinking Water Supply)

Begin Date: 92/08/26 09:15

B/N/Acid Scan *** Continued *** Matrix Spike #1			Water-Total Result Units		
Dibenzofuran	78	%	Recov		
Benzo(ghi)perylene	77	%	Recov		
Indeno(1,2,3-cd)pyrene	71	%	Recov		
Benzo(b)fluoranthene	77	%	Recov		
Fluoranthene	78	%	Recov		
Benzo(k)fluoranthene	75	%	Recov		
Acenaphthylene	87	%	Recov		
Chrysene	72	%	Recov		
Retene	NAF	%	Recov		
4,6-Dinitro-2-methylph+	75J	%	Recov		
1,3-Dichlorobenzene	57	%	Recov		
2,6-Dinitrotoluene	82	%	Recov		
N-Nitroso-di-n-Propyla+	72	%	Recov		
D4-1,2-Dichlorobenzene	40	%	Recov		
4-Chlorophenyl-phenyle+	72	%	Recov		
bis(2-Chloroisopropyl)+	75	%	Recov		
Surrog: 2-Fluorobiphen+	72	%	Recov		
Surrog: 2-Fluorophenol	64	%	Recov		
Surrog: D14-Terphenyl	63	%	Recov		
PYRENE-D10 (SS)	66	%	Recov		
Surrog: D5-Nitrobenzene	69	%	Recov		
Surrog: D5-Phenol	55	%	Recov		
B/N/Acid Scan *** Continued *** Matrix Spike #2			Water-Total Result Units		
Hexachlorobutadiene	67	%	Recov		
Pentachlorophenol	86	%	Recov		
2,4,6-Trichlorophenol	76	%	Recov		
2-Nitroaniline	99	%	Recov		
2-Nitrophenol	106	%	Recov		
Naphthalene, 1-Methyl-	NAF	%	Recov		
Naphthalene	80	%	Recov		
2-Methylnaphthalene	59	%	Recov		
2-Chloronaphthalene	79	%	Recov		
3,3'-Dichlorobenzidine	NAF	%	Recov		
2-Methylphenol	85	%	Recov		
1,2-Dichlorobenzene	70	%	Recov		
o-Chlorophenol (2-Chlo+	90	%	Recov		
2,4,5-Trichlorophenol	89	%	Recov		
Nitrobenzene	89	%	Recov		
3-Nitroaniline	REJ	%	Recov		
4-Nitroaniline	92	%	Recov		
4-Nitrophenol	68	%	Recov		
Benzyl Alcohol	80J	%	Recov		
4-Bromophenyl-phenylet+	84	%	Recov		
2,4-Dimethylphenol	81	%	Recov		
4-Methylphenol	86	%	Recov		
1,4-Dichlorobenzene	69	%	Recov		
4-Chloroaniline	130	%	Recov		
Phenol	73	%	Recov		
bis(2-Chloroethyl)Ether	92	%	Recov		
bis(2-Chloroethoxy)Met+	88	%	Recov		
BIS(2-ETHYLHEXYL) PHTH+	74J	%	Recov		
Di-n-Octyl Phthalate	73J	%	Recov		
Hexachlorobenzene	88	%	Recov		
Anthracene	90	%	Recov		
1,2,4-Trichlorobenzene	68	%	Recov		
2,4-Dichlorophenol	80	%	Recov		
2,4-Dinitrotoluene	92J	%	Recov		
Pyrene	94	%	Recov		
Dimethylphthalate	90	%	Recov		
Dibenzofuran	88	%	Recov		
Benzo(ghi)perylene	87	%	Recov		
Indeno(1,2,3-cd)pyrene	82	%	Recov		
Benzo(b)fluoranthene	88	%	Recov		
Fluoranthene	88	%	Recov		
Benzo(k)fluoranthene	87	%	Recov		
Acenaphthylene	99	%	Recov		
Chrysene	86	%	Recov		
B/N/Acid Scan *** Continued *** Matrix Spike #2			Water-Total Result Units		
Benzo(a)pyrene	81	%	Recov		
2,4-Dinitrophenol	98J	%	Recov		
Dibenzo(a,h)anthracene	82	%	Recov		
Benzo(a)anthracene	85	%	Recov		
4-Chloro-3-Methylphenol	82J	%	Recov		
Benzoic acid	21J	%	Recov		
Hexachloroethane	60	%	Recov		
Hexachlorocyclopentadi+	23	%	Recov		
Isophorone	86	%	Recov		
Acenaphthene	86	%	Recov		
Diethylphthalate	92J	%	Recov		
Di-n-Butylphthalate	86J	%	Recov		
Phenanthrene	90	%	Recov		
Butylbenzylphthalate	80J	%	Recov		
N-Nitrosodiphenylamine	131J	%	Recov		
Fluorene	90	%	Recov		
Carbazole	NAF	%	Recov		

(Sample Complete)



Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Laboratory: EPA, Manchester

Sample No: 92 352354

Description: WHC01

Source: Well (Test/Observation)

Begin Date: 92/08/26 10:20

B/N/Acid Scan			B/N/Acid Scan			Tent Ident - B/N/Aci		
Water-Total			Water-Total			Water-Total		
Result Units			Result Units			Result Units		
+-----+-----+			+-----+-----+			+-----+-----+		
Benzo(a)pyrene	41U	ug/l	Di-n-Octyl Phthalate	41UJ	ug/l	NAPHTHALENE, 1,6,7-TRI+	480NJ*	ug/l
2,4-Dinitrophenol	510UJ	ug/l	Hexachlorobenzene	41U	ug/l			
Dibenzo(a,h)anthracene	100U	ug/l	Anthracene	41U	ug/l			
Benzo(a)anthracene	41U	ug/l	1,2,4-Trichlorobenzene	41U	ug/l			
4-Chloro-3-Methylphenol	200UJ	ug/l	2,4-Dichlorophenol	41U	ug/l			
Benzoic acid	510UJ	ug/l	2,4-Dinitrotoluene	100UJ	ug/l			
Hexachloroethane	41U	ug/l	Pyrene	19J*	ug/l			
Hexachlorocyclopentadi+	200U	ug/l	Dimethylphthalate	41U	ug/l			
Isophorone	41U	ug/l	Dibenzofuran	41U	ug/l			
Acenaphthene	54 *	ug/l	Benzo(ghi)perylene	41U	ug/l			
Diethylphthalate	41UJ	ug/l	Indeno(1,2,3-cd)pyrene	41U	ug/l			
Di-n-Butylphthalate	41UJ	ug/l	Benzo(b)fluoranthene	41U	ug/l			
Phenanthrene	230 *	ug/l	Fluoranthene	15J*	ug/l			
Butylbenzylphthalate	100UJ	ug/l	Benzo(k)fluoranthene	41U	ug/l			
N-Nitrosodiphenylamine	510UJ	ug/l	Acenaphthylene	41U	ug/l			
Fluorene	150 *	ug/l	Chrysene	41U	ug/l			
Carbazole	200UJ	ug/l	Retene	41U	ug/l			
Hexachlorobutadiene	100U	ug/l	4,6-Dinitro-2-methylph+	510UJ	ug/l			
Pentachlorophenol	200U	ug/l	1,3-Dichlorobenzene	41U	ug/l			
2,4,6-Trichlorophenol	100U	ug/l	2,6-Dinitrotoluene	100U	ug/l			
2-Nitroaniline	100U	ug/l	N-Nitroso-di-n-Propyla+	41U	ug/l			
2-Nitrophenol	100U	ug/l	4-Chlorophenyl-phenyle+	41U	ug/l			
Naphthalene, 1-Methyl-	840 *	ug/l	bis(2-Chloroisopropyl)+	41U	ug/l			
Naphthalene	130 *	ug/l	Surrog: 2-Fluorobiphen+	114	% Recov			
2-Methylnaphthalene	630 *	ug/l	Surrog: 2-Fluorophenol	59	% Recov			
2-Chloronaphthalene	41U	ug/l	D4-1,2-Dichlorobenzene	116	% Recov			
3,3'-Dichlorobenzidine	1000U	ug/l	Surrog: D14-Terphenyl	90	% Recov			
2-Methylphenol	41U	ug/l	PYRENE-D10 (SS)	77	% Recov			
1,2-Dichlorobenzene	41U	ug/l	Surrog: D5-Nitrobenzene	0	% Recov			
o-Chlorophenol (2-Chlo+	41U	ug/l	Surrog: D5-Phenol	0	% Recov			
2,4,5-Trichlorophenol	200U	ug/l						
Nitrobenzene	41U	ug/l						
3-Nitroaniline	REJ	ug/l						
4-Nitroaniline	510U	ug/l						
4-Nitrophenol	260U	ug/l						
Benzyl Alcohol	820UJ	ug/l						
4-Bromophenyl-phenylet+	41U	ug/l						
2,4-Dimethylphenol	41U	ug/l						
4-Methylphenol	41U	ug/l						
1,4-Dichlorobenzene	41U	ug/l						
4-Chloroaniline	510U	ug/l						
Phenol	41U	ug/l						
bis(2-Chloroethyl)Ether	41U	ug/l						
bis(2-Chloroethoxy)Met+	41U	ug/l						
BIS(2-ETHYLHEXYL) PHTH+	41UJ	ug/l						

(Sample Complete)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Blank ID: BW2245D1

B/N/Acid Scan			B/N/Acid Scan		
Blank #1			*** Continued ***		
Result Units			Result Units		
Benzo(a)pyrene	8U	ug/l	Di-n-Octyl Phthalate	8UJ	ug/l
2,4-Dinitrophenol	100UJ	ug/l	Hexachlorobenzene	8U	ug/l
Dibenzo(a,h)anthracene	20U	ug/l	Anthracene	8U	ug/l
Benzo(a)anthracene	8U	ug/l	1,2,4-Trichlorobenzene	8U	ug/l
4-Chloro-3-Methylphenol	40UJ	ug/l	2,4-Dichlorophenol	8U	ug/l
Benzoic acid	100UJ	ug/l	2,4-Dinitrotoluene	20UJ	ug/l
Hexachloroethane	8U	ug/l	Pyrene	8U	ug/l
Hexachlorocyclopentadi+	40U	ug/l	Dimethylphthalate	8U	ug/l
Isophorone	8U	ug/l	Dibenzofuran	8U	ug/l
Acenaphthene	8U	ug/l	Benzo(ghi)perylene	8U	ug/l
Diethylphthalate	8UJ	ug/l	Indeno(1,2,3-cd)pyrene	8U	ug/l
Di-n-Butylphthalate	8UJ	ug/l	Benzo(b)fluoranthene	8U	ug/l
Phenanthrene	8U	ug/l	Fluoranthene	8U	ug/l
Butylbenzylphthalate	20UJ	ug/l	Benzo(k)fluoranthene	8U	ug/l
N-Nitrosodiphenylamine	0.6J*	ug/l	Acenaphthylene	8U	ug/l
Fluorene	8U	ug/l	Chrysene	8U	ug/l
Carbazole	40UJ	ug/l	Retene	8U	ug/l
Hexachlorobutadiene	20U	ug/l	4,6-Dinitro-2-methylph+	100UJ	ug/l
Pentachlorophenol	40U	ug/l	1,3-Dichlorobenzene	8U	ug/l
2,4,6-Trichlorophenol	20U	ug/l	2,6-Dinitrotoluene	20U	ug/l
2-Nitroaniline	20U	ug/l	N-Nitroso-di-n-Propyla+	8U	ug/l
2-Nitrophenol	20U	ug/l	4-Chlorophenyl-phenyle+	8U	ug/l
Naphthalene, 1-Methyl-	8U	ug/l	bis(2-Chloroisopropyl)+	8U	ug/l
Naphthalene	8U	ug/l	Surrog: 2-Fluorobiphen+	69	% Recov
2-Methylnaphthalene	8U	ug/l	Surrog: 2-Fluorophenol	81	% Recov
2-Chloronaphthalene	8U	ug/l	D4-1,2-Dichlorobenzene	26	% Recov
3,3'-Dichlorobenzidine	200U	ug/l	Surrog: D14-Terphenyl	83	% Recov
2-Methylphenol	8U	ug/l	PYRENE-D10 (SS)	87	% Recov
1,2-Dichlorobenzene	8U	ug/l	Surrog: D5-Nitrobenzene	82	% Recov
o-Chlorophenol (2-Chlo+	8U	ug/l	Surrog: D5-Phenol	74	% Recov
2,4,5-Trichlorophenol	40U	ug/l			
Nitrobenzene	8U	ug/l			
3-Nitroaniline	REJ	ug/l			
4-Nitroaniline	100U	ug/l			
4-Nitrophenol	50U	ug/l			
Benzyl Alcohol	160UJ	ug/l			
4-Bromophenyl-phenylet+	8U	ug/l			
2,4-Dimethylphenol	8U	ug/l			
4-Methylphenol	8U	ug/l			
1,4-Dichlorobenzene	8U	ug/l			
4-Chloroaniline	100U	ug/l			
Phenol	8U	ug/l			
bis(2-Chloroethyl)Ether	8U	ug/l			
bis(2-Chloroethoxy)Met+	8U	ug/l			
BIS(2-ETHYLHEXYL) PHTH+	8UJ	ug/l			

(Sample Complete)

Project: TEC-575A AVERY RR DUMP AND ROUNDHOUSE

Officer: MZR

Account: FA10PUZZ

Blank ID: BW2245D2

+-----+   B/N/Acid Scan      Water-Total   Blank #2            Result    Units +-----+			+-----+   B/N/Acid Scan      Water-Total                      *** Continued ***   Blank #2            Result    Units +-----+		
Benzo(a)pyrene	8U	ug/l	Di-n-Octyl Phthalate	8UJ	ug/l
2,4-Dinitrophenol	100UJ	ug/l	Hexachlorobenzene	8U	ug/l
Dibenzo(a,h)anthracene	20U	ug/l	Anthracene	8U	ug/l
Benzo(a)anthracene	8U	ug/l	1,2,4-Trichlorobenzene	8U	ug/l
4-Chloro-3-Methylphenol	40U	ug/l	2,4-Dichlorophenol	8U	ug/l
Benzoic acid	100UJ	ug/l	2,4-Dinitrotoluene	20U	ug/l
Hexachloroethane	8U	ug/l	Pyrene	8U	ug/l
Hexachlorocyclopentadi+	40U	ug/l	Dimethylphthalate	8U	ug/l
Isophorone	8U	ug/l	Dibenzofuran	8U	ug/l
Acenaphthene	0.4J*	ug/l	Benzo(ghi)perylene	8U	ug/l
Diethylphthalate	8U	ug/l	Indeno(1,2,3-cd)pyrene	8U	ug/l
Di-n-Butylphthalate	8U	ug/l	Benzo(b)fluoranthene	8U	ug/l
Phenanthrene	1J*	ug/l	Fluoranthene	8U	ug/l
Butylbenzylphthalate	20U	ug/l	Benzo(k)fluoranthene	8U	ug/l
N-Nitrosodiphenylamine	100UJ	ug/l	Acenaphthylene	8U	ug/l
Fluorene	8U	ug/l	Chrysene	8U	ug/l
Carbazole	40UJ	ug/l	Retene	8U	ug/l
Hexachlorobutadiene	20U	ug/l	4,6-Dinitro-2-methylph+	100UJ	ug/l
Pentachlorophenol	40U	ug/l	1,3-Dichlorobenzene	8U	ug/l
2,4,6-Trichlorophenol	20U	ug/l	2,6-Dinitrotoluene	20U	ug/l
2-Nitroaniline	20U	ug/l	N-Nitroso-di-n-Propyla+	8U	ug/l
2-Nitrophenol	20U	ug/l	4-Chlorophenyl-phenyle+	8U	ug/l
Naphthalene, 1-Methyl-	4J*	ug/l	bis(2-Chloroisopropyl)+	8U	ug/l
Naphthalene	0.6J*	ug/l	Surrog: 2-Fluorobiphen+	68	% Recov
2-Methylnaphthalene	3J*	ug/l	Surrog: 2-Fluorophenol	80	% Recov
2-Chloronaphthalene	8U	ug/l	D4-1,2-Dichlorobenzene	29	% Recov
3,3'-Dichlorobenzidine	200U	ug/l	Surrog: D14-Terphenyl	81	% Recov
2-Methylphenol	8U	ug/l	PYRENE-D10 (SS)	84	% Recov
1,2-Dichlorobenzene	8U	ug/l	Surrog: D5-Nitrobenzene	74	% Recov
o-Chlorophenol (2-Chlo+	8U	ug/l	Surrog: D5-Phenol	76	% Recov
2,4,5-Trichlorophenol	40U	ug/l			
Nitrobenzene	8U	ug/l			
3-Nitroaniline	100U	ug/l			
4-Nitroaniline	100U	ug/l			
4-Nitrophenol	50U	ug/l			
Benzyl Alcohol	160U	ug/l			
4-Bromophenyl-phenylet+	8U	ug/l			
2,4-Dimethylphenol	8U	ug/l			
4-Methylphenol	8U	ug/l			
1,4-Dichlorobenzene	8U	ug/l			
4-Chloroaniline	100U	ug/l			
Phenol	8U	ug/l			
bis(2-Chloroethyl)Ether	8U	ug/l			
bis(2-Chloroethoxy)Met+	8U	ug/l			
BIS(2-ETHYLHEXYL) PHTH+	0.9J*	ug/l			

(Sample Complete)